

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033023\
 Data File : BN024643.D
 Acq On : 30 Mar 2023 15:17
 Operator : CG/JU
 Sample : SST0.885
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8247

Quant Time: Mar 30 16:18:45 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 30 15:33:41 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	9646	0.400	ng/u1	0.00
4) Naphthalene-d8	10.817	136	28225	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.638	164	15731	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.384	188	33346	0.400	ng/u1	0.00
17) Chrysene-d12	21.574	240	28540	0.400	ng/u1	0.00
23) Perylene-d12	24.056	264	21182	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	7930	0.776	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.396	152	28026	0.803	ng/u1	0.00
18) Fluoranthene-d10	19.408	212	55884	0.795	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	8650	0.829	ng/u1	95
5) Naphthalene	10.867	128	55529	0.796	ng/u1	99
7) 2-Methylnaphthalene	12.473	142	34754	0.804	ng/u1	100
8) 1-Methylnaphthalene	12.693	142	36337	0.802	ng/u1	99
10) Acenaphthylene	14.360	152	44210	0.803	ng/u1	100
11) Acenaphthene	14.703	153	38882	0.805	ng/u1	100
12) Fluorene	15.684	166	44367	0.809	ng/u1	99
14) Pentachlorophenol	17.033	266	5424	0.868	ng/u1	99
15) Phenanthrene	17.426	178	72811	0.792	ng/u1	100
16) Anthracene	17.519	178	61382	0.822	ng/u1	99
19) Fluoranthene	19.441	202	80503	0.797	ng/u1	100
20) Pyrene	19.803	202	80088	0.772	ng/u1	100
21) Benzo(a)anthracene	21.557	228	68456	0.800	ng/u1	99
22) Chrysene	21.612	228	74191	0.772	ng/u1	100
24) Benzo(b)fluoranthene	23.293	252	70651	0.811	ng/u1	96
25) Benzo(k)fluoranthene	23.345	252	69820	0.845	ng/u1	95
26) Benzo(a)pyrene	23.948	252	54067	0.783	ng/u1	95
27) Indeno(1,2,3-cd)pyrene	26.650	276	65895	0.806	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.670	278	53210	0.819	ng/u1	97
29) Benzo(g,h,i)perylene	27.451	276	56368	0.794	ng/u1	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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